AMENDMENTS TO THE CLAIMS

This listing of the claims will replace all prior versions, and listings, of claims in the application:

LISTING OF THE CLAIMS

1. (currently amended) A compound having the ~~structural~~ formula $L^1[MQ^1Q^2]L^2$ in which M is a ~~mid-transition~~ metal selected from the group consisting of Nb, Ta, Mo, W, Mn and Re, Q^1 and Q^2 are each a univalent ~~radical~~ substituent, and L^1 and L^2 are ligands coordinated to M, wherein each of L^1 and L^2 contains a first coordinating atom that is a nitrogen atom contained within a C=N group, and a second ~~coordinating atom that is either a second nitrogen atom, optionally present in a second C=N group, or an~~ oxygen, sulfur or phosphorus atom.

2. (original) The compound of claim 1, wherein, in each of L^1 and L^2, the second coordinating atom is a second nitrogen atom.

3. (original) The compound of claim 2, wherein, in each of L^1 and L^2, the second nitrogen atom is present in a second C=N group.

4. (original) The compound of claim 3, wherein L^1 and L^2 are identical.

5. (original) The compound of claim 4, wherein the first nitrogen atom in each of L^1 and L^2 is bound to a first substituent R_S, and the second nitrogen atom in each of L^1 and L^2 is bound to a second substituent R_L, wherein the difference in steric bulk between R_S and R_L is sufficient to result in isospecificity when the compound is used as a polymerization catalyst.

6. (original) The compound of claim 3, wherein L^1 and L^2 are different.

7. (original) The compound of claim 6, wherein the first and second nitrogen atoms in the ligand L^1 are bound to a first substituent R_S, and the first and second nitrogen atoms in the ligand L^2 are bound to a second substituent R_L, wherein the difference in steric bulk between R_S and R_L is sufficient to result in syndiospecificity when the compound is used as a polymerization catalyst.

8. (original) The compound of claim 1, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

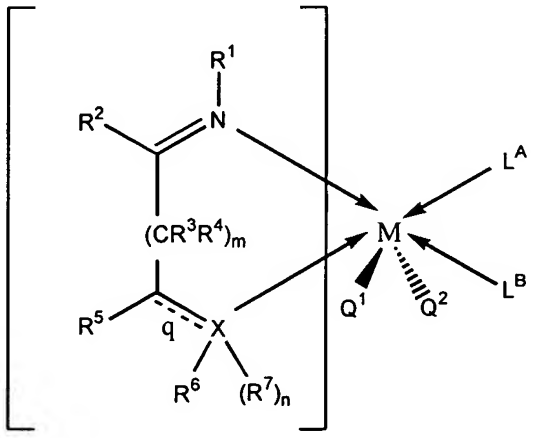
9. (original) The compound of claim 8, wherein y and z are independently integers in the range of 1 to 4 inclusive.

10. (original) The compound of claim 9, wherein y and z are independently 1 or 2.

11. (original) The compound of claim 8, wherein the anions are selected from the group ~~consisting of halide and pseudohalide.~~

12. (currently amended) A compound having the structure of formula (I)

(I)



L^1

wherein:

M is a ~~mid-transition~~ metal selected from Groups VA, VIA and VIIA of the periodic table of the elements;

Q^1 and Q^2 are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted C_1-C_{30} hydrocarbyl, C_1-C_{30} hydrocarbyl substituted with one or more substituents ~~such as electron-withdrawing groups~~, and C_1-C_{30} hydrocarbyl-substituted Group IVB elements, or Q^1 and Q^2 may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group;

m and n are independently zero or 1;

q is an optional double bond;

X is N, O, S or P, with the provisos that (a) when X is N or P, then either n is 1 or q is present as a double bond, but not both, and (b) when X is O or S, then n is zero and q is absent;

R^1, R^6, and R^7 are independently hydrido, hydrocarbyl or substituted hydrocarbyl, and R^2 and R^5 are independently hydrido, halo, hydrocarbyl or substituted hydrocarbyl, or R^1 and R^2 and/or R^5 and R^6 may be taken together to form a linkage -Q-, resulting in a five- or six-membered ring, wherein Q is $-[(CR)_a(Z)_b]-$ in which a is 2, 3 or 4, Z is N, O or S, b is zero or 1, the sum of a and b is 3 or 4, and R is selected from the group consisting of hydrido, halo, hydrocarbyl, hydrocarbyloxy, trialkylsilyl, NR^8_2, OR^9, and NO_2, wherein R^8 and R^9 are each independently hydrocarbyl, or wherein R moieties on adjacent carbon atoms may be linked to form an additional five- or six-membered ring, or R^2 and R^5 may together ~~form a linkage -Q- as just defined~~;

R^3 and R^4 are independently selected from the group consisting of hydrido and hydrocarbyl, or at least one of R^3 and R^4 may be bound through a lower alkylene linkage to an atom contained within L^A or L^B;

L^A and L^B are ligands which may be the same or different and are independently selected from the group consisting of nitrogen-containing, sulfur-containing and oxygen-containing heterocycles, tertiary amines and phosphines, or L^A and L^B may together form a single bidentate ligand that may or may not be the same as L^1,

with the proviso that when (a) L^A and L^B form a single bidentate ligand that is identical to L^1 and M is V or Cr, then either (b) R^1 and R^2 or R^5 and R^6 are taken together to form a linkage -Q- as defined above, or (c) X is other than N, or both (b) and (c).

13. (original) The compound of claim 12, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

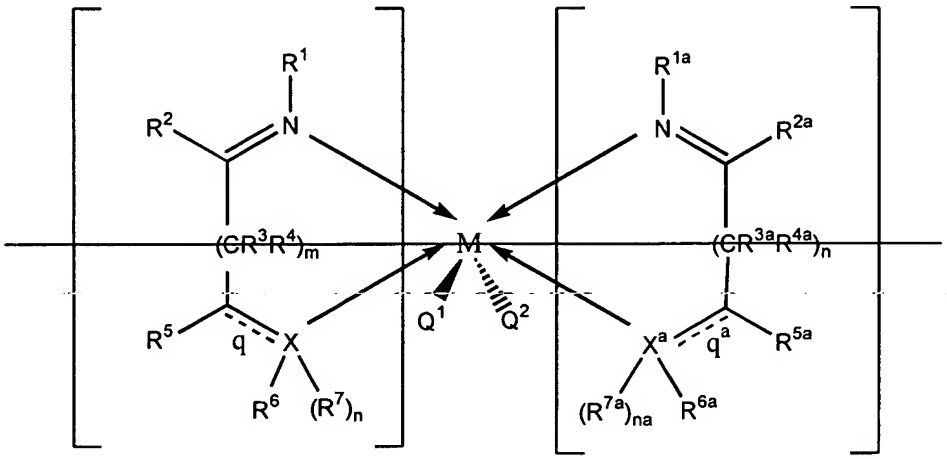
14. (original) The compound of claim 13, wherein y and z are independently integers in the range of 1 to 4 inclusive.

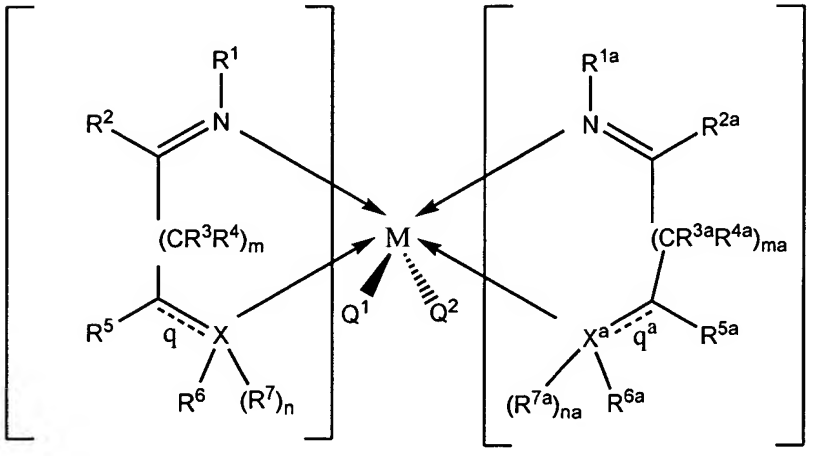
15. (original) The compound of claim 14, wherein y and z are independently 1 or 2.

16. (currently amended) The compound of claim ~~12~~ 13, wherein the anions are selected from the group consisting of halide and pseudohalide.

17. (currently amended) The compound of claim 12, having the structure of formula (II)

(II)





wherein, q^a, ma, na, and R^{1a} through R^{7a} are defined as for q, m, n and R^1 through R^7, respectively.

18. (original) The compound of claim 17, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

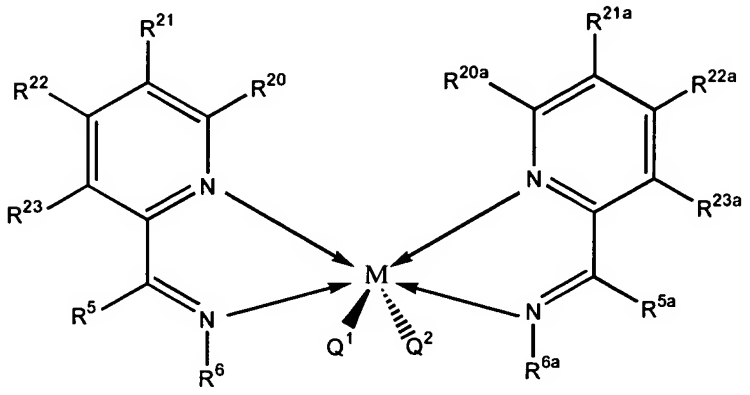
19. (original) The compound of claim 18, wherein y and z are independently integers in the range of 1 to 4 inclusive.

20. (original) The compound of claim 19, wherein y and z are independently 1 or 2.

21. (currently amended) The compound of claim ~~17~~ 18, wherein the anions are selected from the group consisting of halide and pseudohalide.

22. (original) The compound of claim 17, having the structure of formula (V)

(V)



wherein:

R^{20}, R^{20a}, R^{21}, R^{21a}, R^{22}, R^{22a}, R^{23} and R^{23a} are hydrido or hydrocarbyl of 1 to 10 carbon atoms, or any two adjacent R^{20}, R^{20a}, R^{21}, R^{21a}, R^{22}, R^{22a}, R^{23} and R^{23a} groups may be linked to form a five- or six-membered aromatic ring.

23. (original) The compound of claim 22, wherein R^{20}, R^{20a}, R^{21}, R^{21a}, R^{22}, R^{22a}, R^{23} and R^{23a} are hydrido.

24. (original) The compound of claim 22, wherein R^{20} and R^{20a} are methyl, and R^{21}, R^{21a}, R^{22}, R^{22a}, R^{23} and R^{23a} are hydrido.

25. (original) The compound of any one of claims 21, 22 or 23, wherein the compound has a positive charge +y and is associated with y/z anions each bearing a negative charge -z.

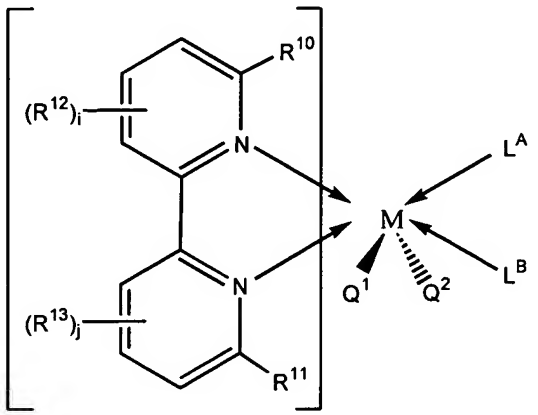
26. (original) The compound of claim 25, wherein y and z are independently integers in the range of 1 to 4 inclusive.

27. (original) The compound of claim 26, wherein y and z are independently 1 or 2.

28. (original) The compound of claim 25, wherein the anions are selected from the group consisting of halide and pseudohalide.

29. (currently amended) A compound having the structure of formula (III)

(III)



L^3

wherein:

M is a ~~mid-transition~~ metal selected from Groups VA, VIA and VIIA of the periodic table of the elements;

Q^1 and Q^2 are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted C_1-C_{30} hydrocarbyl, C_1-C_{30} hydrocarbyl substituted with one or more substituents ~~such as electron-withdrawing groups~~, and C_1-C_{30} hydrocarbyl-substituted Group IVB elements, or Q^1 and Q^2 may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group;

L^A and L^B are ligands which may be the same or different and are independently selected from the group consisting of nitrogen-containing, sulfur-containing and oxygen-containing heterocycles, tertiary amines and phosphines, or L^A and L^B may together form a single bidentate ligand that may or may not be the same as L^3;

i and j are independently zero, 1, 2 or 3; and

R^{10}, R^{11}, R^{12} and R^{13} are independently hydrido, hydrocarbyl or substituted hydrocarbyl.

30. (withdrawn) A catalyst system comprising the compound of any one of claims 1, 5, 7, 12, 17 or 22 and a catalyst activator effective to produce a catalytically active ionic species when combined with said compound.

31. (withdrawn) The catalyst system of claim 30, wherein the catalyst activator is aluminum-containing or boron-containing.

32. (withdrawn) The catalyst system of claim 31, wherein the catalyst activator is aluminum-containing.

33. (withdrawn) The catalyst system of claim 32, wherein the catalyst activator is an organoaluminum compound.

34. (withdrawn) The catalyst system of claim 33, wherein the catalyst activator is an alkyl aluminoxane.

35. (withdrawn) The catalyst system of claim 34, wherein the catalyst activator is methyl aluminoxane.

36. (withdrawn) The catalyst system of claim 31, wherein the catalyst activator is boron-containing.

37. (withdrawn) The catalyst system of claim 36, wherein the catalyst activator is a fluorohydrocarbylboron compound.

38. (withdrawn) The catalyst system of claim 37, wherein the catalyst activator is a fluorinated phenylborate.

39. (withdrawn) The catalyst system of claim 30, further including an inert polymerization diluent.

40. (withdrawn) The catalyst system of claim 39, wherein the diluent is a volatile hydrocarbon solvent.

Claims 41-51 (canceled).

52. (new) The compound of claim 12, wherein Q^1 and Q^2 are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted C_1-C_{30} hydrocarbyl, C_1-C_{30} hydrocarbyl substituted with one or more electron-withdrawing groups, and C_1-C_{30} hydrocarbyl-substituted Group IVB elements, or Q^1 and Q^2 may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group.

53. (new) The compound of claim 29, wherein Q^1 and Q^2 are independently selected from the group consisting of hydrido, halide, alkoxy, amido, unsubstituted C_1-C_{30} hydrocarbyl, C_1-C_{30} hydrocarbyl substituted with one or more electron-withdrawing groups, and C_1-C_{30} hydrocarbyl-substituted Group IVB elements, or Q^1 and Q^2 may together form an alkylidene olefin, acetylene, or a five- or six-membered cyclic hydrocarbyl group.

54. (new) The compound of claim 1, wherein one of L^1 and L^2 contains a second coordinating atom that is either a second nitrogen atom, optionally present in a second C=N group, or an oxygen, sulfur or phosphorus atom, and the other of L^1 and L^2 contains a second coordinating atom that is either a second nitrogen atom, optionally present in a second C=N group, or a sulfur or phosphorus atom.